R5 is H, F, Cl, Br, (C_1-C_6)-alkyl, CF_3, OCF_3, NO_2, CN, O-(C_1-C_6)-alkyl, CO-(C_1-C_6)-alkyl, (C_0-C_6)-alkylene-COOH, (C_0-C_6)-alkylene-COO-(C_1-C_6)-alkyl or SO_2-(C_1-C_6)-alkyl;

A is H, F, Cl, Br, (C_1-C_6)-alkyl, CF_3, OCF_3, NO_2, CN, O-(C_1-C_6)-alkyl, CO-(C_1-C_6)-alkyl, (C_0-C_6)-alkylene-COOH, (C_0-C_6)-alkylene-COO-(C_1-C_6)-alkyl or SO_2-(C_1-C_6)-alkyl;

R7 is H, (C_1-C_6)-alkyl, (C_0-C_6)-alkylene-aryl, O-(C_1-C_6)-alkyl, O-(C_2-C_6)-alkenyl or O-(C_2-C_6)-alkynyl, wherein said (C_1-C_6)-alkyl, (C_0-C_6)-alkylene-aryl, O-(C_1-C_6)-alkyl, O-(C_2-C_6)-alkenyl and O-(C_2-C_6)-alkynyl are optionally mono- or polysubstituted by F, Cl or Br;

R8 is -(C=O)-X;

X is OH, O-(C_1-C_6)-alkyl, NH_2, NH-(C_1-C_6)-alkyl or N-$[(C_1-C_6)$-alkyl$]_2$;

~~m is 1 or 2;~~ and

n is 1 or 2;

and pharmaceutically acceptable salts thereof.

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